

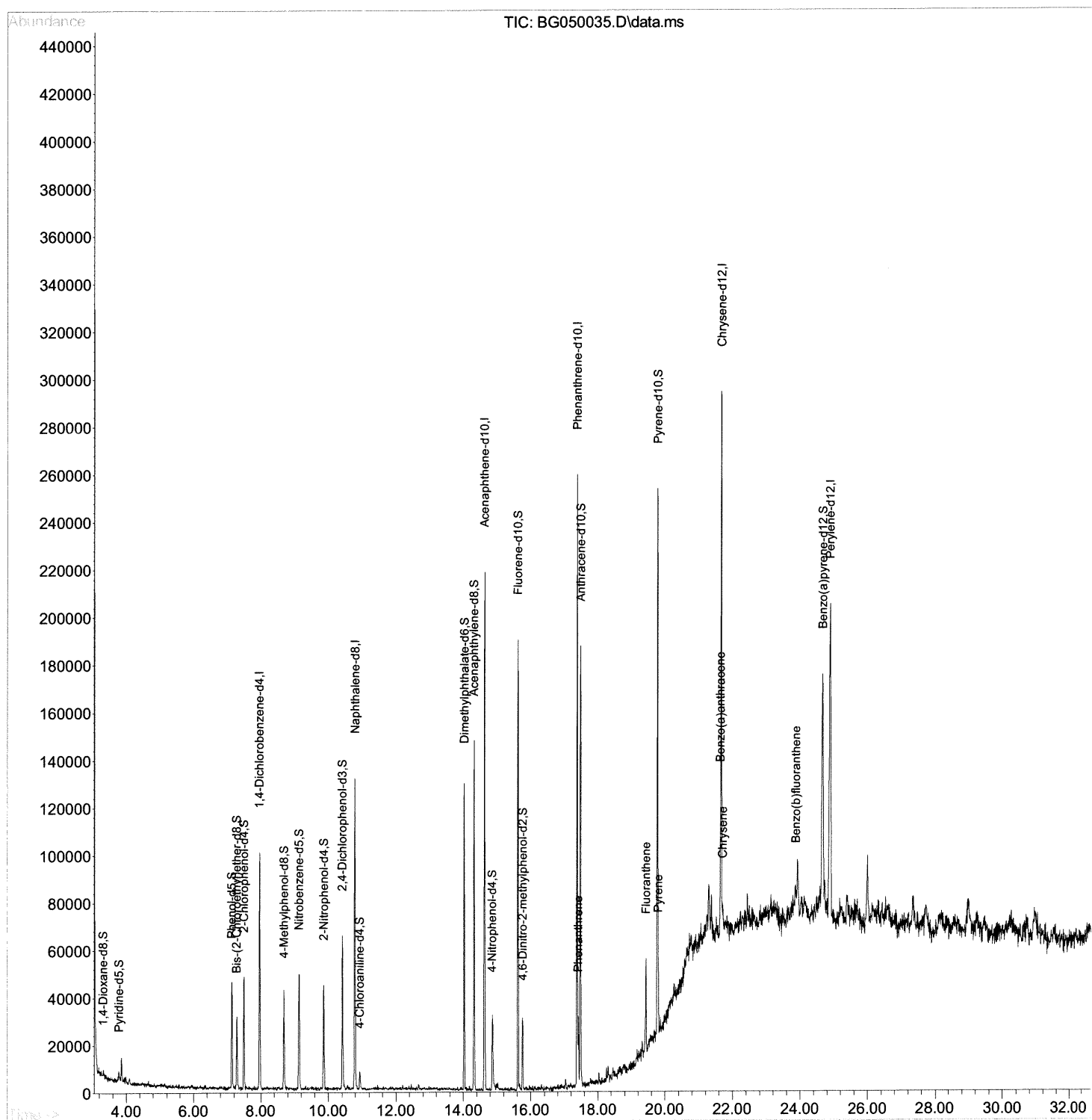
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050035.D
Acq On : 30 Aug 2021 17:04
Operator : CG/JU
Sample : M3458-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AR8

Manual Integrations
APPROVED

mohammad
8/31/2021 11:35:18 AM

Quant Time: Aug 30 17:43:50 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 14:59:55 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

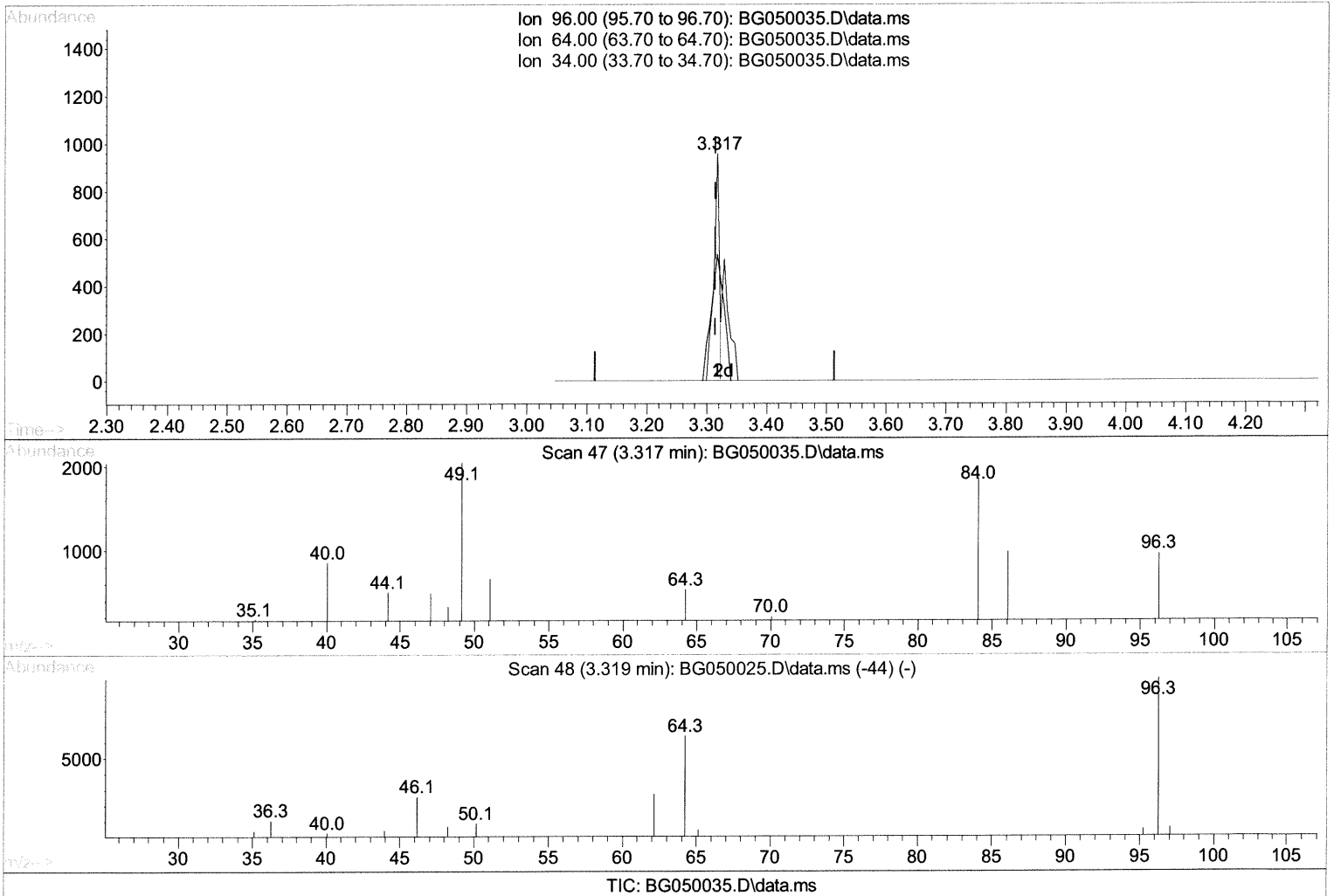
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(3) 1,4-Dioxane-d8 (S)

3.317min (+ 0.003) 1.20 ng/uL

response 626

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	55.65
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

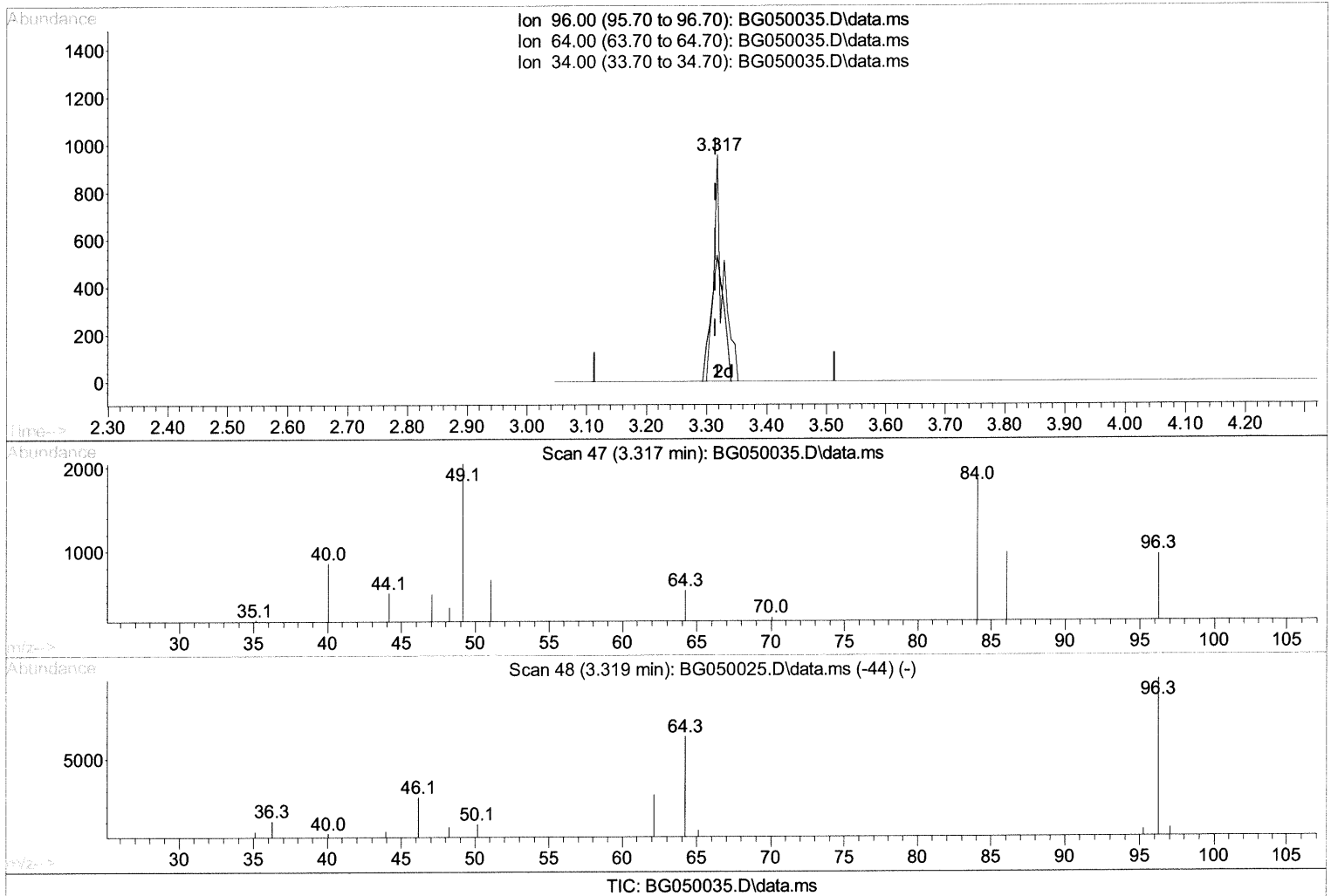
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(3) 1,4-Dioxane-d8 (S)

3.317min (+ 0.003) 1.97 ng/uL m 09/01/21 JU

response 1030

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	55.65
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

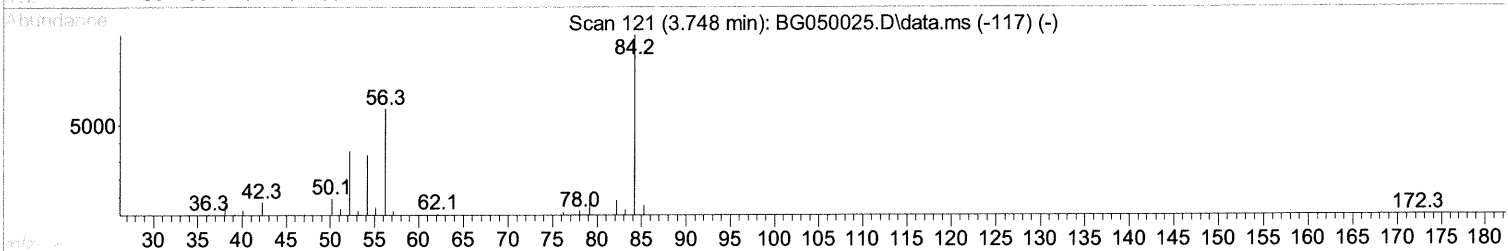
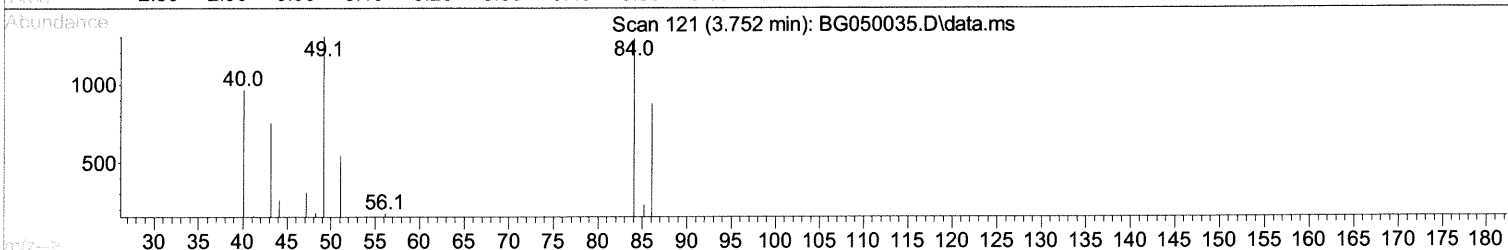
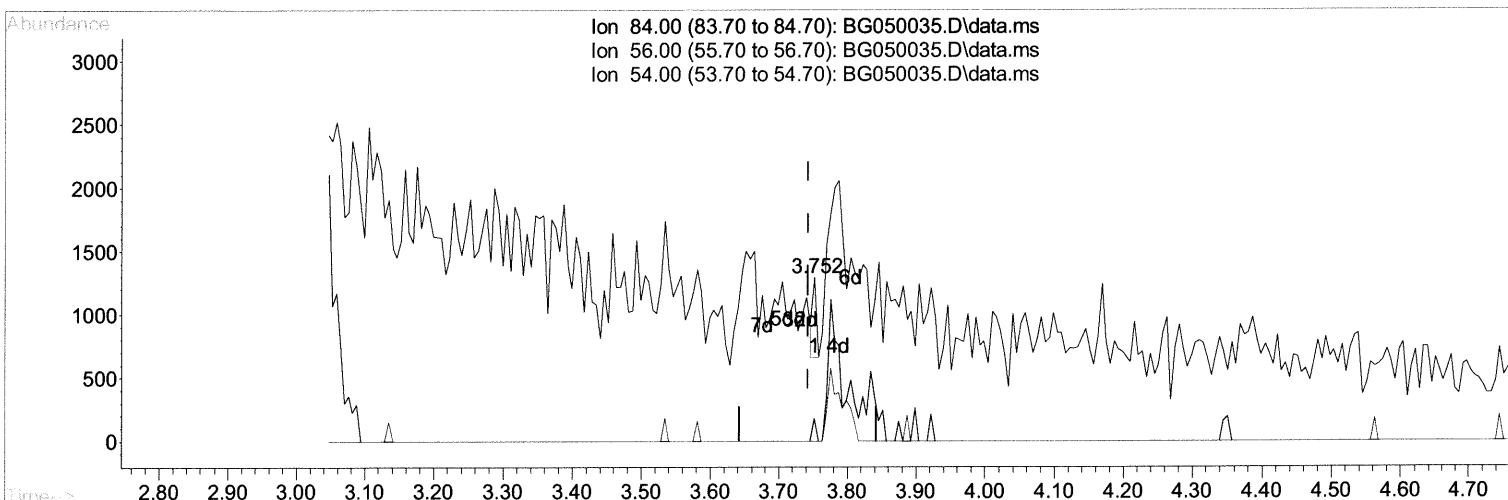
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TIC: BG050035.D\data.ms

(4) Pyridine-d5 (S)

3.752min (+ 0.009) 0.14 ng/ul

response 222

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	13.64#
54.00	37.00	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

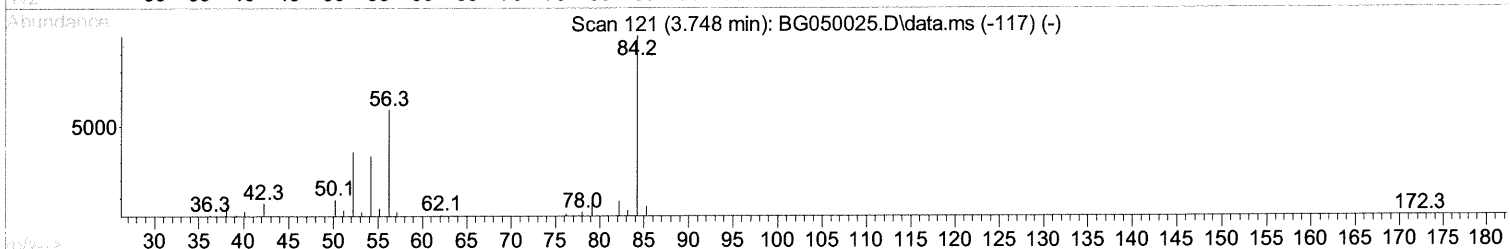
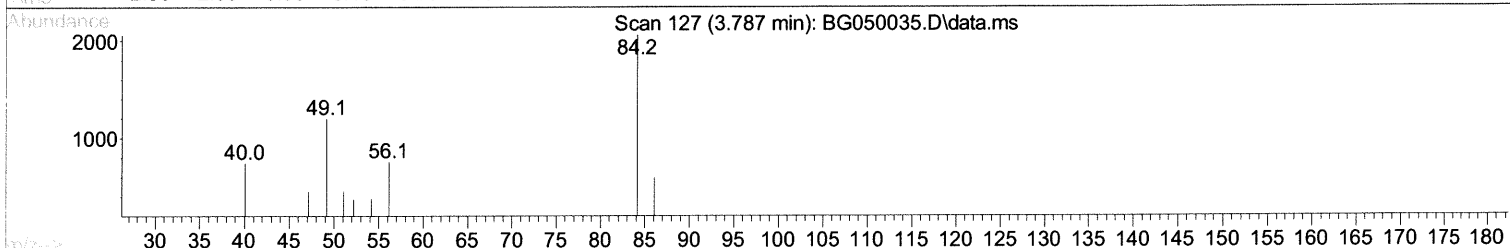
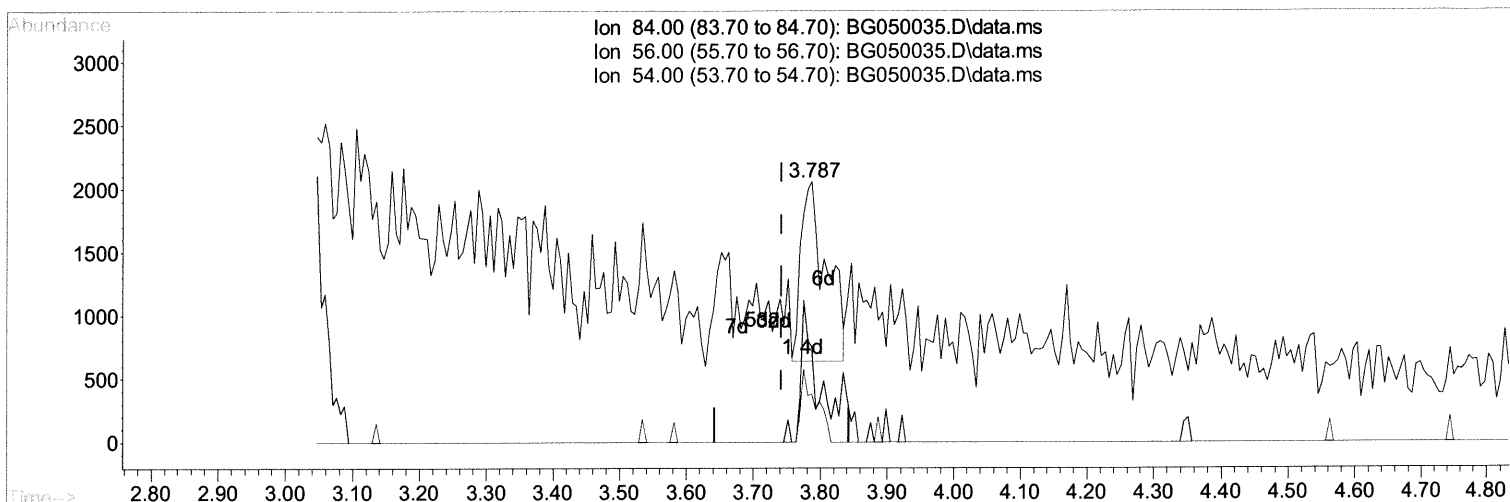
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 Misc :
 ALS Vial : 8 Sample Multiplier: 1

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TIC: BG050035.D\data.ms

(4) Pyridine-d5 (S)

3.787min (+ 0.044) 2.36 ng/ul m 09/01/21JU

response 3724

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	54.30	36.69#
54.00	37.00	18.39#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	27009	20.000	ng/ul	0.00
20) Naphthalene-d8	10.783	136	113545	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.625	164	77407	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.380	188	156277	20.000	ng/ul	0.00
79) Chrysene-d12	21.651	240	131506	20.000	ng/ul	0.00
88) Perylene-d12	24.882	264	136752	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	1030m	1.967	ng/uL	0.00
4) Pyridine-d5	3.787	84	3724m	2.360	ng/ul	0.04
7) Phenol-d5	7.135	99	26912	11.734	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.282	67	15655	12.286	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	21918	12.772	ng/ul	0.00
15) 4-Methylphenol-d8	8.692	113	15429	8.542	ng/ul	0.00
21) Nitrobenzene-d5	9.133	128	12054	13.693	ng/ul	0.00
24) 2-Nitrophenol-d4	9.867	143	13129	12.525	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.413	165	24137	12.910	ng/ul	0.00
31) 4-Chloroaniline-d4	10.930	131	5552	2.182	ng/ul	0.00
46) Dimethylphthalate-d6	14.026	166	86758	14.645	ng/ul	0.00
49) Acenaphthylene-d8	14.320	160	109732	15.798	ng/ul	0.00
54) 4-Nitrophenol-d4	14.860	143	9127	10.649	ng/ul	0.01
60) Fluorene-d10	15.618	176	77722	15.474	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.765	200	7613	7.572	ng/ul	0.00
73) Anthracene-d10	17.480	188	110828	15.793	ng/ul	0.00
81) Pyrene-d10	19.771	212	122785	17.278	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.659	264	106052	14.618	ng/ul	0.00
Target Compounds						
72) Phenanthrene	17.421	178	18301	2.349	ng/ul	94
80) Fluoranthene	19.436	202	23203	2.646	ng/ul#	97
82) Pyrene	19.800	202	18169	2.091	ng/ul#	91
85) Benzo(a)anthracene	21.633	228	11144	1.360	ng/ul	98
87) Chrysene	21.698	228	8320	1.074	ng/ul	93
90) Benzo(b)fluoranthene	23.860	252	13944	1.594	ng/ul#	87

(#) = qualifier out of range (m) = manual integration (+) = signals summed